# COMPUTER AIDED SIMULATION OF THE DYNAMICS OF A MANIPULATOR

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1985

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DEPARTMENT OF MECHANICAL ENGINEERING INDIAN INSTITUTE OF TECHNOLOGY, KANPUR MAY, 1985

# COMPUTER AIDED SIMULATION OF THE DYNAMICS OF A MANIPULATOR

A Thesis Submitted
In Partial Fulfilment of the Requirements
for the Degree of

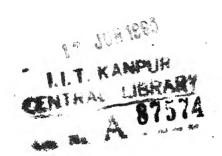
MASTER OF TECHNOLOGY

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by
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to the

DEPARTMENT OF MECHANICAL ENGINEERING
INDIAN INSTITUTE OF TECHNOLOGY, KANPUR
MAY, 1985



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#### CERTIFICATE

Certified that the present work entitled

'Computer Aided Simulation of the Dynamics of a

Manipulater has been carried out by Sri Madhu Shekhar, C.

under the supervision of Dr. B.Sahay, and has not been submitted elsewhere for a degree.

Dr. B.Sahay has gone abroad to Canada for three months, and has authorised me to submit the certificate on his behalf and to look after the viva of Sri Madhu Shekhar.

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May 1985

#### ACKNOWLEDGEMENT

I acknowledge with deep gratitude the invaluable help and constant motivation provided to me by my Thesis Advisor Dr. B. Sahay, without which this work would not have reached completion.

I am also highly indebted to Dr. S.G.Dhande, who readily consented to conduct my oral examination in the absence of Dr. B. Sahay.

Various other people have contributed in various ways towards this project: The Computer Centre, Library, and Mechanical Department staff, and of course my friends.

C. MADHU SHEKHAR

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## LIST OF SYMBOLS

N	Number of links of an Articulate Mechanical System
	(AMS)
L i	ith link of the AMS
$\mathtt{T_{i}^{i}}$	ith joint of the AMS
G <sub>i</sub>	Centre of gravity of L <sub>i</sub>
C <sub>i</sub>	Body fixed coordinate frame of $L_{\hat{1}}$
Oi	Origin of C <sub>i</sub>
a	Link length
α	Link twist
9	Joint angle
u	Joint distance
ρ	Position vector of the centre of gravity of the link
m	Mass of the link
J	Inertia tensor of the link
Α	Transition block matrix
В	Jacobian block matrix
q	Internal coordinate vector
Х	External coordinate vector
$\mathbf{P}^{\gamma}$	Drive vector
F	Force block vector
М	Moment block vector

#### ABSTRACT

Dynamic performance is one of the most significant factors in the design of manipulators particularly for fast and accurate robots recently developed.

Until the end of the last decade, the choice of the manipulator, the actuator and the control system units was a subject of free speculation, frequently based on experience but lacking any systematic method. Hence the need for developing certain criteria and procedures for a systematic design of the manipulator. In the design phase, there is no efficient means except simulation to investigate and evaluate the highly non linear and coupled systems. The aim of this work is to create a software tool for the simulation of all the dynamical values and characteristics of manipulator operations in a particular task execution and thus permit a fast evaluation of a great number of different configurations.

#### CHAPTER I

#### INTRODUCTION

In 1920, the Czech writer Carel Kapek wrote a collective drama entitled R.U.R (Rossum's Universal Robot) and with its translation the word Robot entered the English vocabulary.

For half a century the word Robot appeared on the pages of science fiction stories and novels. Issac Assimov coined the name of the trade, 'Robotics', and provided us Roboticists with an ethic. These terms were later adopted by the scientific and technical community.

The robot may be defined as a class of technical systems which imitate or substitute human locomotion or intellectual function. The Robot Institute of America gives a more precise and technical defination: 'A robot is a reprogrammable multifunctional manipulator designed to move materials, parts, tools, or specialised devices, through variable programmed motions for the performance of a variety of tasks'.

Although the size, structure, function and architecture of robots vary widely, all of them have certain essential features:

- a) Manipulator
- b) Controller
- c) Actuators
- d) Sensors

Robots posses several specific qualities in both a mechanical and a control sense. In the mechanical sense, a feature specific to manipulation robots is that they are open chain active mechanisms, in contrast to conventional mechanisms in which motion is produced primarily by the so called kinematic degrees of freedom. Also, they have a versatile structure, ranging from open to closed, from one to some other kind of boundary conditions. From a control viewpoint, robot systems represent, multi variable and essentially non-linear automatic control systems. A manipulation robot is also an example of dynamically coupled system, and the control task itself is a dynamic task. A further feature is the redundancy reflected in an excess of the degrees of freedom for producing certain functional movements of articulated mechanical system. While it permits a greater mechanical flexibility in task performance, it complicates the control system by introducing optimising procedures for solving the problem of system redundancy. surplus dof are used to satisfy special additional requests.

A list of the uses of robots is impressive. Robots have been used for operation in dangerous environments, in oceanography, space exploration, agriculture, medicines, etc. In industry, their scope is virtually boundless. The development of automatically controlled industrial robots has witnessed three generations: Programmable, Supervisory Controlled, and Artificial Intelligence robots. The boom which we observe today in the sphere of industrial robots is not a mere tribute to vogue. These 'steel collar workers' are now handling, transforming, assembling and dissembling parts, tools and specialised systems. In contrast to hard automation, robots represent 'soft' automation - flexible, neither product, nor operations, nor industry limited, and immune to obsolescence.

The distance separating Carel Kapek from Cybernetics was covered in a quarter of a century. The next twenty five year period built the foundations of what seemed pure fantasy. In this quarter we are witnessing a quantitative accumulation and a qualitative leap in the field of Robotry.

#### LITERATURE REVIEW

Although Robotics is a multidisciplinary field, incorporating the results of numerous scientific and engineering disciplines, two major areas can be identified: areas

involving mechanisms and areas involving control systems.

In this work, we have confined ourselves only to the former:
in particular to the dynamics of manipulators.

The first work in the field of dynamics of spatial mechanisms was published by N.G. Bruyevich [ ] as far back as in 1937.

In 1963, H.J. Fletcher, L. Rongved and E.Y.Yu [ studied the motion of a satellite composed of two rigid bodies, connected by a universal joint, under the load due to gravitation.

In 1965, W.W. Hooker and G. Margulien [2] inspired by the preceeding work, studied the general case, where N+l bodies are connected by means of joints with one or two rotational degrees of freedom. Although this method was a significant advance in that it used matrix formalism, it was not able to obtain the matrices as functions of system state.

R.E. Roberson and J. Wittenburg [ in their approach have defined the system of bodies as a graph, whereby the elaborated and known graph properties are used.

The method of W.W. Hooker [ provides a new possibility of eliminating constraints but presents a serious problem in deriving the motion equations.

The method of P.W. Likins [ proposed in 1971, is a more direct application of the preceeding method. Moreover, without reducing the problem's generality, it simplifies the numerical designation of the links in the chain.

J Wittenburg [4] generalised the method to systems having the structure of a topological branch whose joints permit r revolute dof's and t linear dof's. It does not however, include fully closed chains, nor does it offer an explicit matrix procedure for obtaining the equations.

In 1968, J.J. Uicker [3] proposed a method based on Lagrange's equations for the study of dynamical behaviour of joint connected systems of arbitrary structure. The method is sufficiently general to enable motion equations to be simulated on a computer. To overcome certain deficiencies in this method, M. Renaud [ proposed a method based on matrix calculus. In 1977, J. Zabala [5] proposed an algorithm based on M. Renaud's method for automatically formulating motion equations.

In 1974, E.P. Popov et al [ proposed an algorithm for solving the inverse problem based on the Gauss principle of numerical minimisation. The direct problem is solved from the necessary conditions of minimum.

In 1979, M. Vukobratovic and V. Potkonjak [ ] proposed a method based on Appel's equations, and is computer oriented to a great extent.

All the methods based on the general theorem of dynamics (except the method of P.W. Likins) are based on direct enumeration. However for an efficient computer solution, a recursive formulation is desirable. The methods using Newton's and Euler's methods are in principle complex due to the complexity of eliminating the constraints by forces and moments. Lagrangian equations provide the possibility of directly regarding the equations as functions of system control inputs. However, the inherent unsuitability of applying these equations lies in the need to calculate partial derivatives of Lagrangian functions.

#### OUTLINE OF THE PRESENT WORK

Modelling of the dynamics of Open Chain Active

Mechanisms (OCAM) is a central point in the modern approach
to manipulator design for two main reasons. One is connected
with dynamic control, and the second is the development of
procedures for optimal design of manipulators.

As a rule, the methods for the dynamic analysis of active mechanisms use generalised coordinates. However,

in practice such a solution is insufficient, because one needs to consider the so called functional robot motion. This is a motion satisfying certain practical demands. It is therefore necessary to obtain drives producing these functional motions. Considerations of functional dynamics are closely connected with control and conversly. Hence, the term Dynamic Control, meaning, control based on detailed knowledge of the system dynamic characteristics has been introduced.

action among multiple joints, non linear effects, and vary—
ing inertia depending upon the arm configuration. In prac—
tice until the end of the last decade, the designer proceeded
with the design of each joint mechanism without knowing the
actual characteristics of the multi dof motion. Also, the
choice of the kinematic scheme and its different parameters,
the actuator units, and the control systems was a subject of
free speculation, frequently based on experience but lacking
any system of method. Hence many parameters, and often the
motors were over powered. Hence the need for developing
certain criteria and procedures for a systematic choice of
manipulator configuration. In the design phase there is no
efficient means except simulation to investigate and evaluate
the highly non-linear and coupled systems.

The aim of this work is to develop an interactive program for the simulation of all the dynamical values and characteristics of the manipulator operations in a particular taks execution, permitting the designer to view the characteristics in their entirety, and to evaluate a great number of different configurations as fast as possible.

The program consists of four main parts. The first part defines all the vector, matrix and block matrix operations necessary for the simulation algorithm.

The second part is the simulation algorithm.

The algorithm automatically forms and solves the mathematical model of the robot dynamics using the block matrix method. The inputs for the algorithm are the manipulator configuration, the initial state and the manipulation task. The output of the algorithm are the drives in the joints, the P-N diagram, the energy consumed for the performance of the task, etc.

The third part is the Editor. It is used to perform interactive editing of the task, state or configuration of the manipulator.

The fourth part, the Command Interpreter, intreprets the commands given by the user, demands the data to be provided, if necessary, then selects the desired primitives and executes it. For the prevention of the user's careless errors, the program traps them at various places and issues the relevent messages.

The main program is written in Pascal. The subroutines for the drawing of graphs on the graphics terminal is written in Fortran. For Graphic action, PLOT-10 Interactive Graphics Library (IGL) is used.

#### CHAPTER II

#### MODELLING

#### 2.1 INTRODUCTION:

The algorithms for modelling Open Chain Active Mechanisms (OCAM) can generally be classified into two groups : analytical and computer oriented. Analytical procedures have appeared before Computer Aided (CA) methods and originate within the multibody satellite dynamics. Unfortunately the application of these methods ' by hand ' is very tedious and are subject to mistakes. Also these algorithms include either unsuitable numerical operations or impose some calculations which are not formalised. For these reasons it becomes necessary to develop CA methods for mathematical modelling. It will enable the designer to analyse a number of different configurations and choose the most appropriate one to the future of the device. The development of CA methods which perform real time calculations of robot dynamics is a direct contribution to the synthesis of control algorithm for practical purposes.

Any CA method must satisfy the following requirements:

a) The input data for the algorithm are: robot configuration, robot state, workspace state, and robot task. Using such input data the computer itself

forms and solves the mathematical model, i.e., robot dynamics. In principle three problems of dynamics may be solved: a direct and an inverse one, or a combination of these two.

- b) The algorithm includes no numerical differentiation.

  The present CA methods may be divided into three groups:
  - a) Methods based on the general theorem of dynamics and the Newton-Euler equations,
  - b) Methods based on the second order Lagrange equations.
  - c) Methods based on Gibbs-Appel equations and the Gauss principle.

This program uses the general theorem of dynamics and incorporates the Block Matrix method. The Block Matrix method represents the analytically derived mathematical model but by using suitable block formalism the model reduces to the compact matrix form suitable for solving on a computer.

The aim of each CA method is to derive functions f and g such that

 $\ddot{u} = f(u, \dot{u}, P, mechanism configuration)$ 

 $P = g(u, \dot{u}, \dot{u}, mechanism configuration)$ 

where u is the generalised coordinate vector and P is the drive vector.

The functions f and g are not some explicitly prescribed or derived function. They represent large computation algorighms. The realisation of the algorithms f and g are specific to and characteristic of each method and it depends on the mechanical approach. The algorithms f and g, although mutually inverse, (f represents the direct problem and g the inverse one ) are sometimes realised in rather different ways. Most of the CA methods consider the fifth class kinematic pairs only, i.e., joints permitting only one degree of freedom between the segments. If a compound joint is in question, then it is dissembled into a sequence of fifth class joints with small parameter segments between them.

The following assumptions have been made while forming the mathematical model:

- a) All the links are rigid bodies.
- b) All the articulations are fifth class pairs.
- c) All the articulations are perfect.
- d) There are not kinematic singularities.

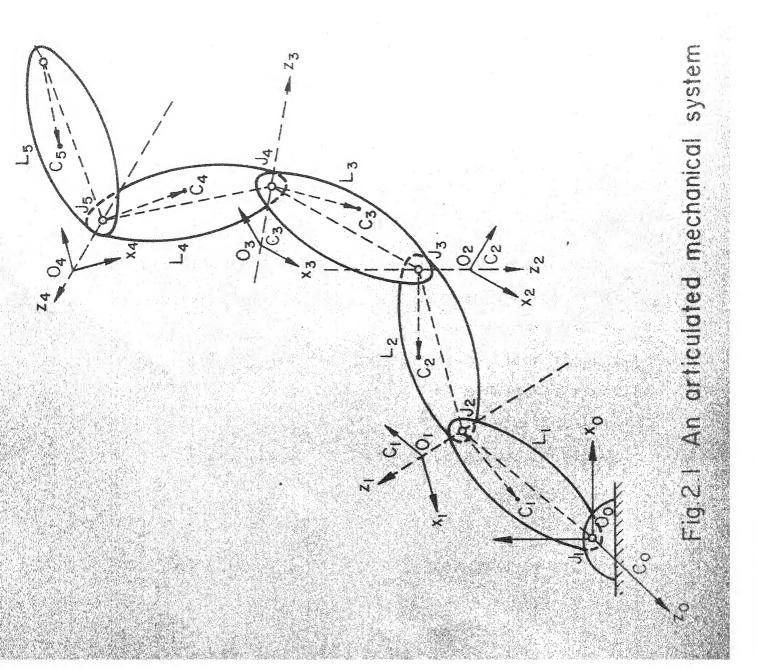
The overall model is developed in three phases: Geometric, Kinematic, and Dynamic.

#### 2.2 THE GEOMETRIC MODEL:

The kinematic structure of the arm has been modelled based on the Denavit and Hartenberg convention [1].

Referring to the Figure 2.1, the links of the Articulated Mechanical System (AMS) are numbered from zero at the base to N at the End Effector (EE). Links  $L_{i-1}$  and  $L_i$  are connected at joint  $T_i$ . The axis of a joint is defined as the axis of rotation for a revolute joint and as a line parallel to the generatrix for a prismatic joint. convenient fixed coordinate system is chosen for the AMS. A Body Fixed (BF) coordinate system  $C_i$  is attached to each link  $L_i$  . The origin of frame  $C_i$  is denoted by  $O_i$ . The z axis of the frame  $C_i$  is oriented along the axis of the joint,  $T_i$  and the x axis is directed along the common normal. The y axis is so chosen as to make a right handed coordinate system. The relative position of successive pair axes is described by the use of the unique common perpendicular between by axes of successive coordinate frames. With each link L; is associated a vector  $\rho_{i}$  defining the centre of mass of  $L_{i}$  with respect to the frame  $C_{i}$ , a mass  $M_{i}$  and an inertia tenser  $J_{i}$ .

Four parameters are used to describe each successive joint and link pair:



Link length, a - distance between successive z axes

Link twist,  $\alpha$  - angle between successive z axes

Joint angle, 0 - angle between successive x axes

Joint distance, u - distance between successive x axes.

Referring to Fig. 2.2, the transformation of the frame  $C_{i-1}$  into the frame  $C_{i}$  may be represented by the concatenation of the following elementary transformations:

a) R ( $z_{i-1}$ ,  $\theta_i$ ) - rotation around  $z_{i-1}$  by an angle  $\theta$  until  $x_{i-1}$  becomes parallel to  $x_i$ 

- b) T  $(z_{i-1}, u_i)$  translation along  $z_{i-1}$  by a distance s until  $x_{i-1}$  coincides with  $x_i$
- c)  $T(x_i, a_i)$  translation along  $x_i$  by a distance a until  $0_{i-1}$  coincides with  $0_i$
- d) R  $(x_i, \alpha_i)$  Rotation around  $x_i$  by an angle  $\alpha$  until  $C_{i-1}$  and  $C_i$  coincide.

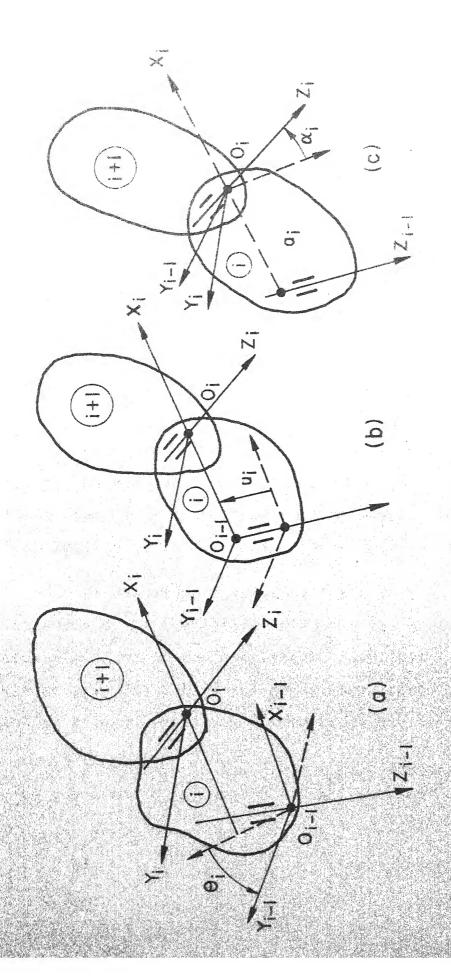


Fig.2.2 Illustrating transition from (i-1)st to ith frame of reference

From these four elementary transformations, a 4 X 4 homogenous transformation matrix  $T_{i-1,i}$ , which relates the position and orientation of frame  $C_i$  to that of frame  $C_{i-1}$  may easily be formed.

$$T_{i-1,i} = \begin{bmatrix} \cos \theta_i & \sin \theta_i & 0 & a_i \cdot \cos \theta_i \\ -\sin \theta_i \cdot \cos \alpha_i & \cos \theta_i \cdot \cos \alpha_i & \sin \alpha_i & a_i \cdot \sin \theta_i \\ \sin \theta_i \cdot \sin \alpha_i & -\cos \theta_i \cdot \sin \alpha_i & \cos \alpha_i & u_i \\ 0 & 0 & 0 & 1 \end{bmatrix} \cdot \cdot \cdot (2.1)$$

Here,  $T_{i-1,i}$  is an orthogonal matrix as it transforms one orthogonal system into another. The transformation between other link coordinate systems may be obtained by the multiplication of intermediate transformation matrices.

It should be noted that for a revolute joint, the joint angle  $\Theta_{\bf i}$  is the joint variable ( the generalised coordinate ), while for a prismatic joint, it is the joint distance  ${\bf u_i}$ . Similarly, the generalised force for a revolute joint is a torque, and for prismatic joint, it is a force.

A point  $r^p$  in the frame  $C_p$  transforms to the point  $r^q$  in the frame  $C_q$  according to the equation

$$\mathbf{r}^{p} = \mathbf{T}_{p,q} \cdot \mathbf{r}^{q} \qquad (2.2)$$

Let  $A_{i-1,i}$  be a 3 X 3 sub matrix of  $T_{i-1,i}$  giving the orientation of the axes of  $C_{i-1}$  with reference to  $C_{i}$ 

$$A_{i-1,i} = \begin{bmatrix} \cos \theta_i & \sin \theta_i & 0 \\ -\sin \theta_i \cdot \cos \alpha_i & \cos \theta_i \cdot \cos \alpha_i & \sin \alpha_i \\ \sin \theta_i \cdot \sin \alpha_i & -\cos \theta_i \cdot \sin \alpha_i & \cos \alpha_i \end{bmatrix} \dots (2.3)$$

And  $l_{i-1,i}$  be 3 X 1 vector giving the position of the origin of the frame  $C_{i-1}$  with reference to  $C_{i}$ 

$$^{1}i-1, i = [a_{i} \cdot \cos \theta_{i} \quad a_{i} \cdot \sin \theta_{i} \quad u_{i}] \qquad \dots (2.4)$$

The equation 2.2 may also be written as

$$r^{i-1} = A_{i-1,i} r^{i} + l_{i-1,i}$$
 ....(2.5)

Then for a vector  $V_{i}$ ,

$$V_{i}^{p} = A_{p,q} V_{i}^{q}$$
 .... (2.6)

where  $A_{p,q}$  is obtained by the multiplication of intermediate transformation matrices

$$r_{i} = \sum_{j=0}^{i} A_{j} l_{j,j} + A_{i} r_{i}^{i} \qquad \dots (2.7)$$

$$V_{i} = \Lambda_{i} V_{i}^{i} \qquad \qquad \dots (2.8)$$

The three columns of the matrix  $A_N$  represents the orientation of the axes of the frame  $C_N$  in terms of the generalised coordinates, and the vector  $\mathbf{l}_N$  represents the position vector of the centre of gravity of the EE in the external coordinate system. Thus the geometric model enables us to convert internal coordinates to external coordinates.

#### 2.3 KINEMATIC MODEL:

The derivation of the equations constituting the kinematic and dynamic model is very lengthy and complex, and is discussed in [6]. The analytical equations themselves are very long. However, by using block matrix formalism, the equations are reduced to a very compact form. The equations, are presented here only in their final form [7].

Let  $a_{\tt i},\ldots a_{\tt N}$  be a set of vectors, and let us define the Block vectors a and  ${\tt a}^{\tt O}$  of dimension ( NX1 (3X1))

$$a = [a_1^1, \dots a_N^N]^T$$

$$a^{\circ} = [a_1 \dots a_N]^T$$

Let  $E_3$  be a 3 X 3 unit matrix, and  $E_N$ , a N X N unit matrix let V be the ( N X N ( 3 X 3 )) block matrix.

$$V = \begin{bmatrix} E_3 & 0 & \cdots & 0 \\ E_3 & E_3 & \cdots & 0 \\ \vdots & \vdots & \vdots & \vdots \\ E_3 & E_3 & \cdots & E_3 \end{bmatrix}$$

Further, let us define the link type matrix, S ( N X N ), and the block matrix  $\boldsymbol{\nu}$  ( N X N (3 X 1))as follows

$$S = \operatorname{diag} \left[ \begin{array}{ccc} s_1 & s_2 & \dots & s_N \end{array} \right]$$

$$v_1 = \operatorname{diag} \left[ \begin{array}{ccc} v_1 & v_2 & \dots & v_N \end{array} \right]$$

$$\dots (2.9)$$

where  $v_i = [O \sin \alpha_i \cos \alpha_i]^T$ 

and s is an indicator, whose value is 0 if the ith joint is revolute and 1 if it is prismatic.

Let A be the orientation block matrix ( N X N  $3 \times 3$ )

$$A = \begin{bmatrix} E_3 & 0 & 0 & \dots & 0 \\ A_{21} & E_3 & 0 & \dots & 0 \\ A_{31} & A_{32} & E_3 & \dots & 0 \end{bmatrix}$$

Now, let us define two functions  $\lambda$  and  $\Lambda$  as follows:

$$\lambda(V_{i}) = \begin{bmatrix} 0 & -V_{iz} & V_{iy} \\ V_{iz} & 0 & -V_{ix} \\ -V_{iy} & V_{ix} & 0 \end{bmatrix}$$
where  $V_{i} = \begin{bmatrix} V_{ix} & V_{iy} & V_{iz} \end{bmatrix}^{T}$ 
...(2.10)

 $A(V) = \text{diag} \left[ \lambda(V_i) \lambda(V_2) \dots \lambda(V_N) \right]$  where  $V_i$  is a vector (3X 1)

Let  $q = [q_1 \ q_2 \ \dots q_N]^T$  be the vector of internal coordinates (N X 1) and  $X = [X_1 \ X_2 \ \dots \ X_N]^T$ , the vector of external coordinates (N X 1). If v is the linear velocity vector, and  $\omega$ , the angular velocity vector, then  $X = [v \ \omega]^T$  is a (2 X 1, (3 X 1)) velocity block matrix.

Now, we have the following relations for velocity block matrix

$$\dot{X} = B\dot{q}$$
 ...(2.12)

where B is a ( 2 X 1( N XN ( 3 X 1))) block matrix

$$B = \begin{bmatrix} B_1 \\ B_2 \end{bmatrix}$$

$$B_1 = -(A\Lambda(1) + \Lambda(\rho) \quad A\nu(E-S) + A\nu S$$

$$B_2 = A\nu(E-S)$$

$$(2.13)$$

For velocities expressed in the external coordinate system, we have

$$\dot{X}^{O} = B^{O} \dot{q}$$
 (2.14)

where

$$B^{\circ} = \begin{bmatrix} B_{1}^{\circ} \\ B_{2}^{\circ} \end{bmatrix}$$
 (2.15)

$$B_{1}^{O} = -\Lambda^{*} (\rho) e^{O} (E-S) \dot{q} + VSe^{O} \dot{q}$$

$$B_{2}^{O} = V e^{O} (E-S) \dot{q}$$

$$\Lambda^{*} (\rho) = \begin{bmatrix} \lambda(\rho_{01}) & 0 & \dots & 0 \\ \lambda(\rho_{02}) & \lambda(\rho_{12}) & \dots & 0 \\ \vdots & & & \ddots \\ \lambda(\rho_{0N}) & \lambda(\rho_{1N}) & \dots & \lambda(\rho_{N-1,N}) \end{bmatrix}$$

Now, let w and  $\epsilon$  be linear and angular acceleration vectors (3 X 1) respectively. Then  $\ddot{X}=[w \ \epsilon]^T$ 

 $\label{eq:weak_problem} \mbox{We have the following relation for acceleration} \\ \mbox{block vector}$ 

$$\ddot{X} = B\ddot{q} + D\dot{q}$$
 ...(2.16)

where

$$D=\begin{bmatrix}D_1\\D_2\end{bmatrix}$$

$$D_{1} = (A \Lambda (1) + \Lambda (\rho)) A \Lambda (\omega) \nu (E-S - (A\Lambda(\omega)\Lambda(1) + \Lambda(\omega) \Lambda(\rho)) A \nu (E-S) + 2A\Lambda(\omega) \nu S \dots$$

$$D_{2} = A (\Lambda(\omega) \nu (E-S) \dots (2.17)$$

In the external coordinate system,

$$\ddot{X}^{O} = B^{O}\ddot{q} + D^{O}\dot{q} \qquad ...(2.18)$$

where

$$D_{i1}^{l} = \begin{bmatrix} 1 & \dots & d_{1i}^{l} & 0 & \dots & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots \\ d_{i-1,i}^{l} & d_{i,1,i}^{l} & 0 & \dots & 0 \\ 0 & \dots & 0 & \dots & 0 \end{bmatrix} \dots (2.19)$$

$$d_{kj}^{l} = \lambda (e_{j-1}) \lambda (e_{k}) \rho_{j-1,i} (1-s_{k}) (1-s_{j})$$

$$\begin{bmatrix} 0 & \dots & d_{12}^{2} \dots & d_{1i}^{2} \dots & 0 & \dots & 0 \\ d_{21}^{2} & 0 & \dots & d_{2i}^{2} & 0 & \dots & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ d_{i1}^{2} & d_{i2}^{2} \dots & 0 & 0 & \dots & \dots & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ d_{kj}^{2} = \lambda (e_{k-1}) e_{j-1} (1-s_{k}) s_{j}$$

$$d_{jk}^{2} = d_{kj}^{2}$$

$$\begin{array}{lll} \mathbf{R}_{\mathbf{j}}^{2} & \mathbf{K}_{\mathbf{i}}^{2} & \mathbf{K}_{\mathbf{i}}^{2} & \mathbf{K}_{\mathbf{j}}^{2} \\ \mathbf{R}_{\mathbf{j}}^{2} & = & \mathbf{d}_{\mathbf{k}\mathbf{j}}^{2} \\ \mathbf{D}_{\mathbf{j}}^{0} & = & \left[ \mathbf{D}_{\mathbf{1}}^{2} & \dots & \mathbf{D}_{\mathbf{N}}^{2} \right]^{T} & \text{dimension (NXN (3X1))} \\ \mathbf{D}_{\mathbf{i}}^{2} & = & \mathbf{\hat{q}} & \mathbf{H}_{\mathbf{i}}^{2} \\ \mathbf{\hat{q}} & = & \left[ \mathbf{\hat{q}}_{\mathbf{1}} & \mathbf{E}_{\mathbf{3}} & \mathbf{\hat{q}}_{\mathbf{2}} \mathbf{E}_{\mathbf{3}} & \dots & \mathbf{\hat{q}}_{\mathbf{N}} & \mathbf{E}_{\mathbf{3}} \right] \end{array}$$

We now have the expressions for the velocity and acceleration of the links in the external coordinate system as well as the BF system. These expressions constitute the kinematic model.

#### 2.4 DYNAMIC MODEL:

The dynamic equations are derived on the basis of D ' Alambert's principle by interrupting the chain fictitiously at a joint and balancing the forces and moments. In this way, N scalar equations are obtained, and transformed into matrix form using block formalism.

Let us introduce the following block matrices of dimension (NXN (3X3)).

$$m = \text{diag } [ m_1 E_3 \dots m_N E_3 ]$$

$$J^o = \text{diag } [ J^o_1 \dots J^o_N ]$$

$$A_o = \text{diag } [ [ A_{1o} \dots A_{No} ]$$

The expressions for the block vectors of the resultant inertial forces and moments can be written together as

$$\begin{bmatrix} F_{I}^{\circ} \\ M_{I}^{\circ} \end{bmatrix} = - \frac{1}{2} B^{\circ} q - C^{\circ} q \qquad \dots (2.20)$$

where

$$\mathcal{T} = \begin{bmatrix} m & 0 \\ 0 & J^{\circ} \end{bmatrix}$$

$$\mathcal{C}^{\circ} = \begin{bmatrix} m & D_{1}^{\circ} \\ \Omega^{\circ} J^{\circ} B_{2}^{\circ} + J^{\circ} D_{2}^{\circ} \end{bmatrix}$$

$$\Omega = \Lambda (\omega^{\circ}) \qquad \dots (2.21)$$

Let us further introduce the vector of the drives  ${\tt P}$  of dimension (  ${\tt NX1}$ )

$$B = [P_1 \dots P_N]^T$$

Then the dynamic equations can be written as

$$W\ddot{q} = B'\dot{q} + C'M_E^O + D'(G^O + F_E^O) + P \dots (2.22)$$

where

$$W = B^{\circ} \mathcal{T} B^{\circ}$$

$$B' = B^{\circ} C^{\circ}$$

$$C' = B_{2}^{\circ}$$

$$D' = B_{1}^{\circ}$$

$$\dots(2.23)$$

by introducing

$$U(q,\dot{q}) = B^{\dagger}\dot{q} + C^{\dagger}M_{E}^{O} + D^{\dagger}(G^{O}+F_{E}^{O}) \dots (2.24)$$

the equation acquires the from

$$W(q) \dot{q} = U (q, \dot{q}) + P$$
 ...(2.25)

Thus we have described the dynamics of the AMS by means of N differential equations in matrix form.

#### CHAPTER III

#### SIMULATION

#### 3.1 INTRODUCTION:

The notion of the simulation of dynamics usually involves the solution of the inverse problem of dynamics, i.e. the determination of the motion for prescribed generalised forces. In this case, the simulation is considered somewhat more liberally so that it also includes the notion of the simulation of the direct problem.

The Fig. 3.1 shows the different models and the different levels of control of a typical manipulator organised in a hierarchial fashion. The highest level defines the task to be carried out, the strategic level divides the imposed task into elementary functions, the tactical level performs the distribution of an elementary movement to the motion of each degree of freedom of the robot, and the executive level executes the imposed motion of each dof. The model developed in the previous chapter is used as the basis for the development of the algorithm for the simulation of the manipulator dynamics, and thus the synthesis of the tactical level.

Before proceeding further, it will be useful to note an important point. The task of a manipulator may

Fig. 3.1 The Different Control Levels And Model of a Typical Manipulator.

be defined as the process of transferring the system state from one bounded region of initial states into another bounded region in the state space within a finite settling time, the system state being in the bounded region of the state space during the transfer. The state of the Articulated Mechanical System ( AMS), at any instant may be described by means of a vector of coordinates. The coordinates are called internal coordinates if the elements of the vector represent the position of the articulation at that time ins-If they represent the coordinates of a predefined point on the End Effector (EE) with reference to the external coordinate system, then they are called the external coordinates . However, in practice, the task is prescribed neither by means of internal coordinates, nor the external coordinates, but by the so called functional coordinates, as it is necessary to consider the functional robot motion. This is a motion satisfying certain practical demands. For instance, for a particular operation, the task of positioning the minimal configuration may be given in terms of spherical coordinates, and the task of orientation of the the EE may given in terms of three Eular angles. For another operation it may be more convenient to give the orientation of the EE in terms of a direction and an angle of rotation round it.

Transformation of the functional coordinates into external coordinates is fairly simple, but the transformation of the external coordinates to internal coordinates is extremely complicated not only because the transformation is not explicit, but because it cannot even be numerically approximated due to the complexity of the system. Also, a set of external coordinates may not have a unique set of corresponding internal coordinates. The external coordinates define the position and orientation of the EE, which we may call 'pose' of the manipulator. On the other hand the internal coordinates determine the posture of the manipulator, i.e., the geometrical configuration adopted by the manipulator while holding an object in a particular pose. The posture of a manipulator completely defines it pose, but for a particular pose a menipulator may be in different postures. A 3 dof manipulator has only one posture in general, and a 6 dof one is believed to have 32 postures in general. A human arm can adopt an infinite number of postures.

### 3.2 THE SIMULATION ALGORITHM:

Let us designate by  $\eta$  a function which transforms the internal coordinates into external coordinates.

$$X = \eta (q)$$
 ..... (3.1)

Now, for the operation of the Computer Aided (CA) method of forming the mathematical model, it is necessary to know q, q and q at each time instant. However, only q appears as input since q and  $\dot{q}$  are calculated by integration starting with known initial state  $q^0$  and  $\dot{q}^0$ . So in order to realise the simulation, it is necessary to develop a procedure for calculating q from the known state q,  $\dot{q}$ , and known external coordinate vector, X.

By differentiating equation 3.1 twice,

$$\dot{X} = \frac{d\eta}{dq} \cdot \dot{q} \qquad ...(3.2)$$

$$X = \frac{d\eta}{dq} \cdot \dot{q} + \frac{d^2\eta}{dq^2} \cdot \dot{q}^2 \qquad ...(3.3)$$

$$A = \frac{d\eta}{dq} \cdot \dot{q} + A \qquad ...(3.4)$$

$$X = B \dot{q} + A$$

B is called the Jacobian matrix (order NXN ).

B and A are functions of the state q, q, and are calculated numerically by means of the kinematic model developed in the previous chapter. Then q is calculated by the relation

$$\ddot{q} = B (\ddot{X} - A)$$
 ... (3.5)

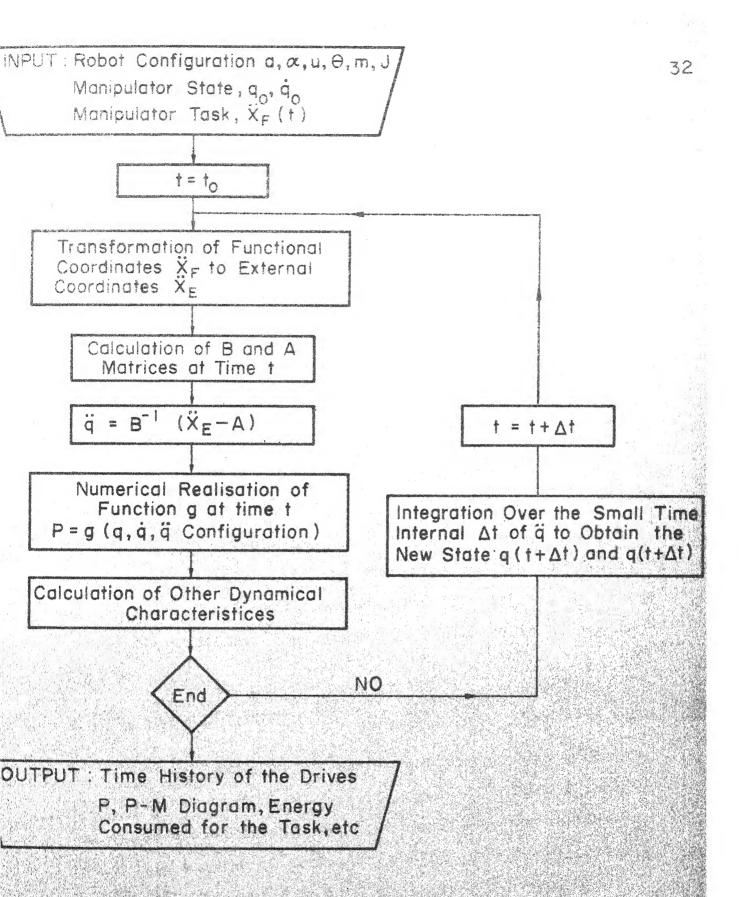


FIG. 3.2 FLOW CHART OF THE SIMULATION ALGORITM

$$\ddot{X}$$
 can be represented by  $\ddot{X} = \begin{bmatrix} W \\ \varepsilon \end{bmatrix}$ 

where w is the linear acceleration, and  $\epsilon$  is the angular acceleration of the centre of gravity of the EE.

From Eqn. 2.16, we can obtain

$$\ddot{X} = B\ddot{q} + D\dot{q}$$

$$= B\ddot{q} + A \qquad ...(3.6)$$

Now from Eqn. 2.25 we can calculate the drives

$$P = g (q, q, q, configuration)$$

$$= Wq-U \qquad ... (3.7)$$

By integrating  $\dot{q}$ , we get the state q,  $\dot{q}$  at the next time instant. With the new state and the corresponding value of the functional coordinates acceleration,  $X_F$  we can repeat the process. The simulation algorithm is presented in Fig. 3.2.

#### 3.3 ADJUSTMENT BLOCKS:

As pointed out earlier, although the manipulation task can be defined fully by means of external coordinates, they are unsuitable for setting. For the various individual classes of tasks, the most suitable variables are given as inputs, which naturally reflect the type of the

task. These are called the, functional coordinates. The transformation of the functional coordinates to the external coordinates is done in a block, known as the adjustment block, within the simulation algorithm.

Before proceeding with the various classes of tasks, let us define some notions precisely. By positioning we mean moving the centre of gravity of the last segment (or some predefined point on it ) to some desired point in the work space according to a prescribed motion law.

Full orientation of the body in space means an exactly determined angular position of the body with respect to the external space.

Partial orientation of the body means that the given body axis coincides with a prescribed direction is space.

To solve the positioning task, which is part of every manipulation task, three dof are necessary.

To solve the positioning task along with the task of partial orientation, five dof are necessary.

To solve the positioning task along with the task of full orientation, six dof are needed.

Now some typical classes of tasks and ways of using the dof will be analysed.

- A) A manipulator with four dof solves the positioning task by using three dof, and with the one remaining performs operations frequently sufficient for many practical tasks.
  - 1. The task is that of positioning in the cartesian coordinate system, i.e. x(t), y(t),z(t) for the centre of gravity of the EE, and the fourth dof is prescribed directly ( $q_{\Delta}(t)$ ).
  - 2. Same as l., but positioning is given in cylindrical coordinates (t),  $\Theta(t)$ , z(t).
  - 3. Same as 1., but positioning is given in spherical coordinates r(t),  $\theta(t)$ ,  $\phi(t)$ .
- B) A manipulator with five dof solves the positioning and partial orientation task.
  - The task is given in the form of positioning (Al , A2 or A3) and partial orientation.
- C) A manipulator with six dof solves the positioning and full orientation task, as well as all the problems in which fewer dof are needed ( when compensation of the dof surplus is done ).

- 1. The task is given in the form of positioning (Al,A2, or A3), and full orientation of the EE in terms of the three Euler angles,  $\Theta(t)$ ,  $\psi(t)$ ,  $\beta(t)$ .
- 2. The task is given in the form of positioning (Al,A2, or A3 ) and full orientation is given in terms of one direction  $\Theta(t)$  and  $\beta(t)$ , and an angle of rotation around it,  $\psi$  (t).
- 3. The task is given in terms of positioning as in 2 plus partial orientation. Motion along one dof is prescribed directly ( $q_k(t)$ ).

The derivation of the adjustment blocks for these classes of task will now be given briefly. We will use the same of matrix notation for  $X_{\rm E}$  (external coordinates)i.e.

$$\dot{X}_{E} = \begin{bmatrix} w \\ \varepsilon \end{bmatrix} = B\dot{q} + A$$
Let  $B = \begin{bmatrix} \Omega \\ r \end{bmatrix}$  and  $A = \begin{bmatrix} \Omega \\ \varphi \end{bmatrix}$ 

 $\overline{\text{Al}}$  The input to the block is  $X_F$  (functional coordinates)

$$\ddot{X}_{F} = \begin{bmatrix} \ddot{x} & \ddot{y} & \ddot{z} & \ddot{q}_{4} \end{bmatrix}^{T}$$
Hence,  $w = \begin{bmatrix} \ddot{x} & \ddot{y} & \ddot{z} \end{bmatrix}^{T}$ 
...(3.7)

Using the equation 3.5  $\ddot{q}_1, \ddot{q}_2$  and  $\ddot{q}_3$  are calculated  $\ddot{q}_4$  is prescribed directly.

$$\underline{A2}$$
  $\ddot{X}_F = [\rho \dot{\theta} \dot{z} \dot{q}_A]^T$ 

The following equations can easily be derived using elementary transformation.

$$w_{x} = w_{\rho} \cos \theta - w_{\theta} \sin \theta$$

$$w_{y} = w_{\rho} \sin \theta + w_{\theta} \cos \theta$$

$$w_{z} = z$$

$$(3.8)$$

where  $w_{\rho} = p - \rho \dot{\theta}^2$ ,  $w_{\theta} = \rho \dot{0} + 2 \dot{\rho} \dot{\theta}$ ,  $w_{z} = \dot{z}$ 

 $\rho$ ,  $\tilde{\rho}$ .  $\Theta$ ,  $\tilde{\Theta}$  etc are obtained by integration of  $\rho$ ,  $\tilde{\Theta}$  etc. the rest is as in Al.

$$\underline{A3}$$
  $\ddot{x}_F = [\ddot{r}, \ddot{\theta}, \dot{\beta}, \dot{q}_A]^T$ .

Again, the following equations can easily be derived. Please refer to Fig. 3.3. a,b,c.

$$w_x = w_r \cos \cos \theta - w_\theta \sin \theta - w_\beta \sin \cos \theta$$
 $w_y = w_r \cos \sin \theta + w_\theta \cos \theta - w_\beta \sin \sin \theta$ 
 $w_z = w_r \sin + w_\beta \cos \theta$ 

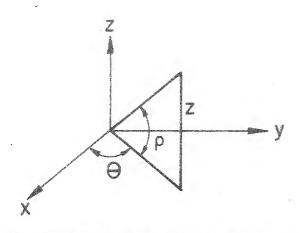
...(3.9)

where

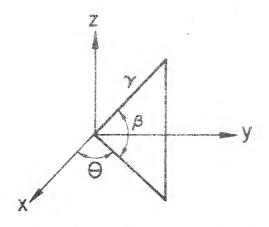
$$w_{\mathbf{r}} = \ddot{\mathbf{r}} - \mathbf{r} \dot{\beta}^{2} - \mathbf{r} \cos^{2} \beta \dot{\theta}^{2}$$

$$w_{\mathbf{\theta}} = 2\dot{\mathbf{r}}\dot{\theta} \cos \beta + \dot{\mathbf{r}}\dot{\theta} \cos \beta - 2\dot{\mathbf{r}}\dot{\beta}\dot{\theta} \sin \beta$$

$$w_{\mathbf{\theta}} = 2\dot{\mathbf{r}}\dot{\beta} + \dot{\mathbf{r}}\dot{\beta} + \mathbf{r} \sin \beta \cos \beta \dot{\theta}$$



a) Cylindrical co-ordinates



(b) Spherical co-ordinates

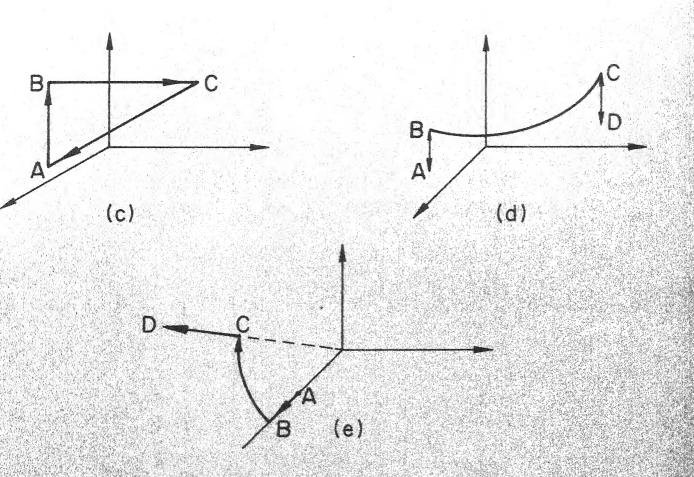


Fig. 3.3 Various trajectories of the manipulator tip

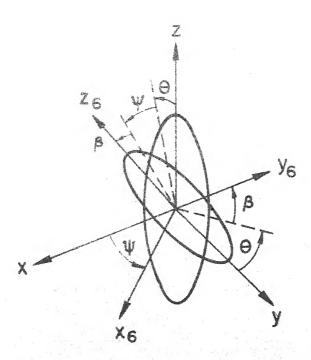


Fig.3.4 Euler angles of the BF system  $(x_6,y_6,z_6)$  relative to the external system (x,y,z)

all the

The positioning may also be prescribed in cylindrical or spherical coordinates. In any case, w is calculated as in Al, A2, or A3, as required. Please refer to Fig. 3.4.

The angular acceleration is given by

$$\varepsilon = A_6 \left( \pi \begin{bmatrix} \ddot{\varphi} \\ \ddot{\psi} \end{bmatrix} + 1_{\pi} \begin{bmatrix} \dot{\psi} & \dot{\xi} \\ \dot{\beta} & \dot{\varphi} \end{bmatrix} \right)$$

where

$$\pi = \begin{bmatrix} \cos \psi & 0 & 1 \\ \sin \psi \cdot \sin \beta & \cos \beta & 0 \\ \sin \psi \cdot \cos \beta - \sin \beta & 0 \end{bmatrix}$$

$$1_{\pi} = \begin{bmatrix} 0 & 0 & -\sin \psi \\ -\sin \beta & \sin \psi \cdot \cos \beta & \cos \psi \cdot \sin \psi \\ -\cos \beta & -\sin \psi \cdot \sin \beta & \cos \psi \cdot \cos \beta \end{bmatrix} \dots (3.10)$$

 $A_{6}$  is the transition matrix of the last segment

$$A_6 = A_1 A_2 \cdots A_6$$

C2 This method is suitable in many practical cases, for eg. spraying powder along a prescribed path, or screwing in a bolt.

$$\ddot{\mathbf{x}}_{F} = \begin{bmatrix} \ddot{\mathbf{x}} & \ddot{\mathbf{y}} & \ddot{\mathbf{z}} & \ddot{\mathbf{\theta}} & \mathbf{\beta} & \ddot{\boldsymbol{\psi}} \end{bmatrix}^{T}$$

w is determined as in Cl

Referring to Fig. 3.5, the direction of any line may be prescribed by means of two angles  $\theta$  and  $\beta$  . Rotation around this line is given by  $\psi$  .

 $\epsilon$  'is calculated from the relation

$$\lambda(\epsilon) = \tilde{A} A^{-1} - [\lambda(\omega)]^2 \qquad \dots (3.11)$$

where

$$A = A_{1} A_{2} A_{3}$$

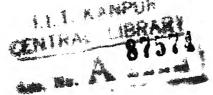
$$A_{1} = \begin{bmatrix} \cos \theta & -\sin \theta & 0 \\ \sin \theta & \cos \theta & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

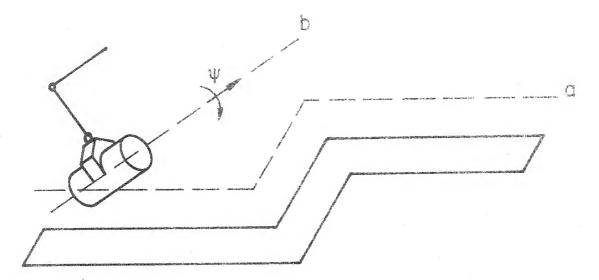
$$A_{2} = \begin{bmatrix} \cos \beta & 0 & -\sin \beta \\ 0 & 1 & 0 \\ \sin \beta & 0 & \cos \beta \end{bmatrix}$$

$$A_{3} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & \cos \psi & -\sin \psi \\ 0 & \sin \psi & \cos \psi \end{bmatrix}$$

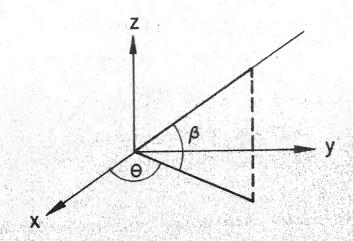
$$A_{4} = A_{1} A_{2} A_{3} + A_{1} A_{2} A_{3} + A_{1} A_{2} A_{3}$$

$$A_{5} = A_{1} A_{2} A_{3} + A_{1} A_{2} A$$





(a) Spraying powder along a prescribed trajectory



(b) External angles  $\Theta$ ,  $\beta$ 

Fig. 3.5

Bl In this case, the manipulator has five dof, for which positioning is prescribed as in Al, A2 or A3 and the orientation is prescribed by the requirement that the EE should coincide with a given direction in space

$$\ddot{X}_F = \begin{bmatrix} \ddot{x} & \ddot{y} & \ddot{z} & \ddot{0} & \beta \end{bmatrix}$$

where the angles  $\Theta$  and specify the direction, as in C2. w is calculated as in C2.  $\epsilon$  is claculated as follows.

$$\varepsilon = \begin{bmatrix} \dot{h}_1 \\ \dot{h}_2 \end{bmatrix}$$

where

$$\dot{h}_1 = -\cos\beta \dot{\beta}^2 \cos\theta - \sin\beta \dot{\beta} \cos\theta + 2 \sin\beta \dot{\beta} \sin\theta \dot{\theta} - \cos\beta \cos\theta \dot{\theta}^2$$

$$-\cos\beta \sin\theta \dot{\theta}$$

$$\dot{h}_2 = -\cos\beta \dot{\beta}^2 \quad \sin\Theta - \sin\beta \dot{\beta} \sin\Theta - - 
 -2\sin\beta \dot{\beta} \cos\Theta \dot{\Theta} - \cos\beta \sin\Theta \dot{\Theta} 
 + \cos\beta \cos\Theta \dot{\Theta} . 
 (3.13)$$

 $\underline{\text{C3}}$  In this case five functional coordinates are prescribed exactly as in Bl and the remaining dof is prescribed directly namely  $\textbf{q}_k$  (t)

$$\ddot{X}_F = [\ddot{x} \ \ddot{y} \ \ddot{z} \ \ddot{\theta} \ \ddot{\beta} \ \ddot{q}_6]$$

The adjustment block is derived as in Bl.

# 3:4 OTHER DYNAMICAL CHARACTERISTICS :

# (a) P N Diagram:

Since the drives and generalised velocities, q, are calculated at each step of the simulation, it is possible to draw a diagram for each joint, the characteristics of the driving motor torque P versus the motor rpm N, i.e. the diagram P-N for each motor. Such a diagram is very useful during the synthesis and choice of the servosystems. The producer gives the  $P_{\rm max}$  - N motor characteristics in the catalogue where  $P_{\rm max}$  is the maximal motor torque at motor rpm N. By comparing the necessary characteristics obtained by means of simulation with the one from the catalogue, one can decide whether the chosen motor suits its applications.

The diagram P-N is obtained in such a way that for each time instant and for joint k,

$$N = R\dot{q}_k$$

$$P = \eta(R) P_{k} / N$$
 ..... (3.14)

where R is the reduction ratio of the subject joint and  $\eta$  (R) is its mechanical efficiency. Repeating the procedure for each time instant, the desired diagram is obtained.

# (b) Energy Consumed:

During the simulation, energy consumed during each step  $\Delta t_{\bf i}$  is found, and the total energy is calculated by summation of these values.

$$E^{i} = E^{i-1} + E_{\Delta t_{i}} \qquad \dots (3.15)$$

where E<sup>i</sup> is the total energy consumed including during the ith time step, and E  $_{\Delta t_i}$  is the energy consumed during the ith time interval. To calculate E  $_{\Delta t_i}$  the everage drive value on the interval is taken

$$P_{\text{med}}^{i} = \frac{1}{2} (P^{i-1} + P^{i})$$
 ....(3.16)

$$E_{\Delta t_{i}} = \Delta q^{i} \cdot P_{med}^{i} \qquad \dots (3.17)$$

where  $\Delta q^i = q^i - q^{i-l}$ . The elements of the vectors  $q^i$  and  $P^i_{med}$  are absolute value. The average drive value is used to avoid more complex interpolation.

We can thus obtain the total energy consumed for the given manipulation task.

Thus, we now have a simulation algorithm which has the following inputs: the manipulator configuration, the manipulator state, and the manipulation task in the form most suited for prescribing ( according to the class of

functional movements). A for the output, we obtain time history of the generalised coordinates, velocities of the drives, the torque-rpm diagram for each actuator, and the total energy consumed.

## CHAPTER 1.V

#### THE PROGRAM MODULES

## 4.1 INTRODUCTION:

The simulation algorithm described in the previous chapter is implemented on a DEC 1090 mainframe in an interactive, user friendly environment.

In one session, the user can set the input data, run the simulation routine and obtain the results on a VDU, a graphic terminal or the line printer, and save the output on the disk for future reference. If not satisfied with the results, the user may change any of the input parameters on line using the Editor routine, and rerun the Simulator routine. To effect the above actions is the Command Interpreter routine, which accepts commands from the user, decodes it, and performs the necessary action. The commands are arranged in a hierarchial fashion with three levels. Errors, if any, resulting either from user action or during the program run are trapped, and relevant warnings or messages given to the user.

## 4.2 THE PROGRAM MODULES:

The program consist of four main modules.

These will now be discussed in detail.

## 4.2.1 Block Matrix Opera or :

This module consists of a number of procedures which define all the vector, matrix and block matrix operations necessary for the simulation algorithm. Let us use the following notation to describe the vectors, matrices, and block matrices:

VEC-3D A vector of dimension 1X3

VEC-ND A vector of dimension 1XN

MAT-3D A matrix of dimension 3X3

MAT-ND A matrix of dimension NXN

BLK-V-V A block vector of dimension 1XN, whose elements are of type VEC-3D

BLK-V-M A block vector (1XN) with elements of type
MAT-3D

ELK-M-V A block ma rix (NXN) with elements of type VEC-3D

BLK-M-M A block matrix (MXN) with elements of types MAT-3D

BLK-V-V A block matrix (1X2) with elements of type

BLK-V-V

BLK-V-M-V A block matrix (1X2) with elements of type BLK -M-V

BLK-M-M-M A block matrix (2X2) with elements of type BLK-M-M.

An operation is named based on the type of the output variable. For instance, if a procedure takes two entities of type MAT-ND and after performing operations returns an entity of type MAT-ND, it will be named as OPN-MND. Simularly, if a procedure returns an entity of type BLK-M-M, it will be named as OPN-M-M. However, for operations involving dot product of vectors, the keyword DPN will be used instead of OPN. For describing the operation, the following notation will be used:

The following is the list of all the operations defined in this module. The last entity is always the output parameter.

i) OPN-V3D ( OPCODE: integer; S: real; A: MAT-3D; B,C: VEC-3D; D: VEC-3D);

Here and in all further operations, OPCODE represents the operation code, whose value determines which operation is to be performed on the input parameters. In this case, the following operations are performed.

OPCODE	OPERATION
1	D = B+C
2	D = B-C
3	$D = S_{\bullet}B$
4	$D = A \cdot B$

If the OPCODE is 1, then the vectors B and C are added according to the rules of the vector addition. Dummy variables may be passed for S and A. In this program, for dummy entities, a zero, or a null vector, or a null matrix, or a null block matrix is passed, depending upon the operations.

- ii) OPN- VND ( OPCODE: integer; S:real; A,B,C: VEC-ND); In Pascal, unlike Fortran, all the variables used must be declared, and the type checking is done at compile time. Hence, OPN-V3D is seperated from OPN-VND, although the two differ only in the dimensions of the parameters. The OPCODE and the operations are exactly same as for i).

This is for doing operations on matrices. The Operations for the OPCODES are as follows

OPCODE	OPERATION
1	C = A + B
2	C = A - B
3	$C = A^T$
4	$C = S \cdot A$
5	$C = A \cdot \cdot B$

This procedure is used, for instance, to calculate the transition matrix  $A_{\rm N}$  in Eqn. 2.6 by repeatedly applying it. Further, it is used, along with OPN-V3D and OPN-VND, in almost all the operation that follow.

This is same as OPN-M3D, except for the dimensions of the matrices A,B, and C, and the OPCODES have the same meaning. This procedure is called while evaluating the expressions in Eqn. 2.12 through 2.18. It is also called by some of the procedure that follow.

v) OPN1-V-V (OPCODE: integer; S:VEC-ND; A:BLK-M-V; B,C,D: BLK-V-V);

The operations performed by this procedure for different values of OPCODE is as follows

OPCODE	OPERATION
ĺ	D = B + C
2	D = B - C
3	$D = S \cdot A$

It is easy to see that while performing the operation D=B+C, the elements of B are added to the corresponding elements of C. These elements are not numbers, but vectors. Hence, to calculate D, this procedure invokes OPN-V3D, N times. Also, in this case, dummy variables are passed as S and A.

vi) OPN2-V-V (A:BLK-V-M; B:BLK-M-V; C:BLK-V-V);

No OPCODE is given here as only one operation is to be performed, i.e.,

$$C = A \cdot B$$

An instance of the use of this operation is in Eqn. 2.19 for calculating  $\textbf{D}_{i}^{l}$  and  $\textbf{D}_{i}^{2}$  .

vii) OPN-M-V (OPCODE: integer; S: MAT-ND; A: BLK-M-M;
B,C,D:BLK-M-V);

The OPCODE and corresponding operations are as follows;

OPCODE		OPERATION
	*	D = B + C
0		D = B - C
3		$D = B^{T}$
4		$D = S \cdot B$
5		$D = A \cdot B$

This procedure calls OPN-V3D repeatedly to evaluate D. Also, when OPCODE is 3, the transpose of B is evaluated, not the elements of B. This procedure is used in Eqn. 2.15 through 2.19 to evaluate the expressions.

This operates on block matrices whose elements are matrices, and is used for evaluating the block matrices B, D and C in Eqn. 2.13 through 2.18 and 2.21. The OPCODES and the corresponding operations are as follows:

OPCODE	OPERATIONS
1	C = A + B
2	C = A - B
3	$C = A^{T}$
4	$C = S \cdot A$
5	$C = A \cdot B$

Again, it should be noted that to evaluate C, the elements of A and the corresponding ones in B are operated by OPN-M3D repeatedly as the elements are not numbers but matrices.

ix ) OPN1-V-V-V (S:VEC-ND; A:BLK-V-M-V; B:BLK-V-V-V); As there is only one operation to be performed, OPCODE is not used. This does the operation

$$B = A \cdot S$$

To evaluate expressions like Bq and Dq in Eqn. 2.12 and 2.16, this procedure is used.

x) OPN2-V-V (A,B,C: BLK-V-V-V);

This does the operation

$$C = A + B$$

Please note that as the elements of A and B are block matrices of type BLK-V-V, OPN-V-V is invoked repeatedly to evaluate C = A + B. This is used in Eqn. 2.16.

xi) OPN-V-M-V (A:BLK-V-M-V; B: BLK-M-M-M; C:BLK-V-M-V); This is used in Eqn.2.23 and does the operation

 $G = B \cdot A$ 

xii) DPN-R (A,B: VEC-3D; R: real);

This gives R as the dot product of two vectors A and B, and is used in the procedures that follow.

xiii) DPN-V (A: BLK-M-V;B:BLK-V-V; C:VEC-ND);
This performs the operation

$$C = B \cdot A$$

To do this, DPN-R is called repeatedly as the elements of B and A are vectors.

xiv) DPN1-M (A,B: BLK-M-V; C: MAT-ND);

This does the operation

 $C = A \cdot B$ 

xv) DPN2-M (A,B: BLK-V-M-V; C:MAT-ND);
This does the operation

# $C = A \cdot B$

by repeatedly applying DPN1-N to the elements of A and the corresponding ones of B and adding the matrices so obtained by OPN-MND.

There is also a procedure for calculating the inverse of a matrix.

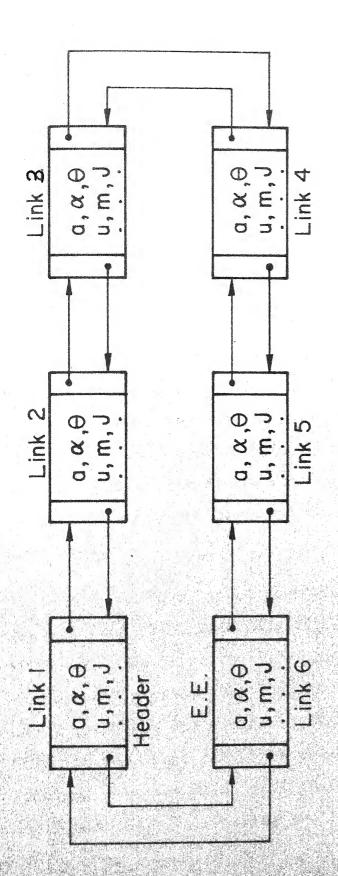
xvi) INVERSE (DIM: integer; A, B: MAT-ND);

B is returned as the inverse of A. DIM is the dimension of the matrices A and B. This procedure is used to calculate the inverse of the Jacobian matrix, in Eqn. 3.5

#### 4.2.2 Simulator:

This module implements the simulation algorithm described in the previous chapter. This again consists of seven main submodules.

The first submodule creates the link structures for storing the input data and initialises certain variables and files. The configurations parameters of the links of the manipulator are represented internally by means of a doubly linked circular list, and are accessed by means of



The link structure for representation of the manipulator configuration Fig. 4.-

a pointer, as shown in Fig. 4.1. One node of the link is designated as the Header, and contains a record consisting of all the configuration parameters like linklength, link twist, mass, inertia, centre of gravity, etc. Also, the record contains two pointers, a left pointer pointing to the last link, and a right pointer pointing to the second In the same way, the left pointer of the second link contains the address of the pointer to the first link, and the right pointer, of the third. When the last link is reached its right pointer is linked to the first link. In the same manner, the generalised joint variable, q, velocity, d and the transition matrix A are stored in another link The acceleration, in functional coordinates, describing the task of the manipulator, is stored as an array ( for a series of time instants ) of an array ( for each dof).

reading from files TASK.DAT, CONFIG.DAT, CMND.DAT and PROG.HLP. The first two files obviously are for task and configuration data, The sets of commands used by the Command Interpreter are stored in CMND.DAT, and PROG.HLP contains help for the user, which can be accessed while the program is running. The output files POSE.DAT, POSTR.DAT, DRIVE.DAT and PNDIAG.DAT, for storing the data for internal coordinates,

external coordinates, drives, and the P-N diagram respectively, are also initialised. Global variables like unit matrices, unit block matrices, matrices E and V in Eqn. 2.13 and 2.15, null matrices, etc. are initialised. Certain option are also set to their default values. The system of units for both input and output data are set to MKS units, the representation for positioning and orientation are set to Cartesion and Euler respectively (Section 3.3)

The second part of this module forms and solves the geometric and the kinematic model. First, the functions  $\lambda$  and  $\Lambda(\text{Eqn. 2.10})$  are defined in procedures LAMDA and BIG-LAMDA. Another function, LAMDASTAR is defined, which takes two indices i and j, and returns the matrix  $\lambda(\rho_{i,j})$  (Eqn. 2.15). Next comes the procedure KINEMATIC. The geometric and kinematic modelling is done here. First the transition matrices are computed in the body fixed system, (Eqn.2.3) and then the matrices so obtained are concatenated to obtain the transition matrices in the external coordinate system. Next the block matrix B<sup>O</sup> is evaluated (Eqn.2.15). Now, the block matrix D<sup>O</sup> is evaluated (Eqn. 2.19). Finally, internal velocities and accelerations are computed (Eqn. 2.14, 2.18).

The third submodule computes the Jacobian matrix and the vector A from the matrices  $B^{O}$  and  $D^{O}$  evaluated earlier. (Eqn. 3.4).

The fourth submodule is the adjustment block. The acceleration in functional coordinates are transformed into those in external coordinates, as discussed in Section 3.3.

Using the Jacobian matrix and the vector A obtained in submodule three, the accelerations in external coordinates obtained in submodule four are transformed into those in internal coordinates, in the submodule five.

The sixth submodule is for dynamic modelling. Eqn. 2.20 through 2.25 are used and the drives, at a time t, obtained. Also, the torque and rpm are calculated for each joint at this time instant, as well as the energy consumed since the time instant ( $t-\Delta t$ ), from the Eqn. 3.14 and 3.17.

In the seventh submodule, the new state, q and  $\dot{q}$  are computed by the integration of  $\dot{q}$  over the time interval  $\Delta t$  , and the time is incremented by  $\Delta t$  .

The first two main modules have now been discussed. But the user will really not be concerned with these two modules. To use the program, what he is interested in is the commands to be given, and the Editor.

# 4.2.3 Command Interpreter (CI):

This is for performing simulation, editing, input/output, etc., as and when requested by the user. The commands are organised in a hierarchial fashion into three levels. The commands can be shortened until it becomes unique. If the command is misspelt, or an illegal command or an ambiguous command is given at any level, a message is given to the user, warning this. If a valid command is given then it is interpreted, and the control passed to the module required to perform the action desired.

At the top level, a double arrow'>>' is given as a prompt . Now the user can type in any of the following commands:

Command	Action
>> HELP	Give help for top level actions
>> TYPE INPUT	Display the input data on the
	terminal
>> LIST INPUT	List the input data on the line
	pinter
>> SIMULATE	Perform simulation -
>> TYPE OUTPUT	Display the output data on the
	terminal
>> LIST OUTPUT	List the output data on the line
	pinter

>> PLOT Plot the graphs on the graphics terminal.

>> EDIT Edit the input data.

>> UPDATE Update the input file.

>> DOCUMENT Type the Documentation.

>> EXIT Exit from the top level.

When any of the input, output or edit action are requested by the user, the CI in invoked recursively, and the user enters the second level. For input action, the second level prompt is a vertical line '|'. Now the user can type in any of the following commands:

HELP Type help for the input action commands.

SYSTEM Type the system of units, coordinates,

etc.

CONFIG Type the configuration parameters of

the manipulator.

STATE Type the state parameters of the

manipulator.

TASK Type the task of the manipulator .

EXIT Exit this level.

For output action, the second level prompt is a forward slash '/'. The user may now type in any of the following commands:

/	HELP	Type help for output action
		commands.
/	DRIVE	Display the drives in the joints.
/	ORIENT	Display the joint orientation .
/	POSITION	Display the joint positions .
/	VELOCITY	Display the joint velocities.
/	ACCELER	Display the joint accelerations.
/	PM DIAG	Display the P-M diagram.
/	ENERGY	Display the energy consumed.
/	EXIT	Exit this level.

For PLOT action, the second level prompt is a back slash '  $\backslash$  ' . The commands are same as for output actions.

For performing editing of the input data, when the EDIT command is given, then the user enters the second level. The prompt is a percent sign'%'. Now the following commands can be given:

%	HELP	Type help for edit action.
%	SYSTEM	Change system of units, represen-
		tation of coordinates etc.
%	CONFIG	Edit configuration parameters.
%	STATE	Edit state parameters.
%	TASK	Edit manipulator task.
%	EXIT	End edit.

When one of the SYSTEM, CONFIG, STATE, or TASK command is given, the CI is once again called recursively, and the user thus inters the third level. The prompt is an asterix '\*'. If the user typed the SYSTEM command, the Editor asks for the system parameter to be changed. The user can now give the following answers:

\* System Parameter : UNITS

\* System Parameter : ANGLE

\* System Parameter : POSITION

\* System Parameter : ORIENT

The command interpreter is now once again invoked to determine the system of UNITS (CGS,MKS or FPS), ANGLE (DEGREES or RADIANS), POSITION (CARTESIAN, CYLINDRICAL, or SPHERICAL), ORIENT (NONE, PARTIAL, FULL). If the system of orientation is partial or full, CI is again called to determine whether the functional coordinates are Euler angles, a vector and an angle, etc (as discussed in section 3.3). For instance,

\* System of representation for Orientation: EULER.

In response to the CONFIG command, the editor gives the following prompt,

\* Add node, delete node, or change node?

In response, the user can respond by typing ADD, DELETE, or CHANGE. If the CHANGE command is given, the editor asks for the number of the link whose link parameters are to be changed.

#### \* Link Number:

The user can now typing a number ( l to N ) for the link whose link parameters are to be changed . Now the C.I. is called to determine the particular parameter to be changed. The following commands can be given:

- \* Link parameter : LINKTYPE
- \* Link parameter : LINKLENGTH
- \* Link parameter : LINKTWIST
- \* Link parameter : JOINTDIST
- \* Link parameter : COG
- \* Link parameter : MASS
- \* Link parameter : INERTIA
- \* Link parameter : REDUCT
- \* Link parameter : EFF

After this, the user can type in the new values, and the corresponding changes will be made.

If the DELETE command is given, then the Editor asks for the number of the link to be deleted, and deletes that link. Similarly, for an ADD command, a number K is taken, and a new link added after the Kth link.

If the state parameters are to be edited, then the second level Editor command STATE can be given.

#### % STATE

Now the C.I. is invoked (third level) to determine the state parameter to be changed.

### \* State Parameter:

Now the following commands may be given :

- \* State parameter : INITIAL VELOCITY
- \* State parameter : INITIAL POSIT

Now the vectors may be typed in as directed by the editor. Similarly, by issuing the second level; TASK command, the task parameters may be changed. The editor takes a number K from the user and changes the acceleration vector at the Kth time instant.

The user may make as many changes as the desires, then exit the editor module. Now the top level command SIMULATE may be given to rerun the simulation module. Now, the output parameters may be observed and evaluated. If the user is satisfied with the results, then he may issue the UPDATE command, which will update the input files CNFG.DAT and TASK.DAT. This cycle the user may continue as many times as he desires.

All the valid commands to the CI are stored in the file CMND.DAT as sets of commands. When the CI is called, a parameter is passed to it indicating which set of commands it should accept from the user. In a command set, for each command coreesponds an action to be performed. The action for a particular command may include a call to the CI. By making such recursive calls to the CI, a hierarchial command structure is built. Control can pass to the higher level only after the lower level releases the CI. Such a structure allows new commands to be easily incorporated at any level, to cater to new modules which may be added to the program.

# 4.2.4 Editor Module:

The editor may be called from the top level to interactively edit any of the input parameters. After editing the new input parameters may be saved by updating the file. All the Edit mode commands are described in Section 4.2.3.

Apart from the above four main modules the program also contains two other small modules: The HELP module for giving help to the user at various levels, and the UTILITY module which contains utility routines for performing certain actions which, though not directly related to the actual running of the program, help to make it more useable.

There are standard procedures for reading integers and reals from a device. Suppose the program is waiting for the user to give an real number from the terminal: If by mistake the user types an integer, or some character, the program is interupted, and control returned to the monitor. To avoid this, two procedures, READIN, and READRL are written, which dont read the numbers directly, but as a sequence of characters which are converted into their ASCII code and then the number computed.

The whole program, except for the plotting routine is written in Pascal. The plotting routine, which calls certain routines from the Interactive, Graphics Library, PLOT-10, is written in Fortran, as these routines are only Fortran-Callable. The object code for the plotting routine and the rest of the program is linked together using the Fortran I/O Link Facility, FTNLNK, available here on DEC 10.

# 4.3 SETTING INPUT DATA:

The data regarding the configuration, state, or task of the manipulator are initially given through disk files. These parameters may later be edited from the terminal.

The data for the configuration of the manipulator is given in the file CNFG . DAT in the following format:

NO OF LINKS

GRAVITY FORCE VECTOR

LINKTYPE/LINK LENGTH/LINK TWIST /JOINT ANGLE or JOINT DISTANCE

COG VECTOR

MASS

INERTIA TENSOR

EXTERNAL FORCE VECTOR

EXTERNAL MOMENT VECTOR

REDUCTION RATIO/ JOINT EFFICIENCY

A blank line or comment

LINKTYPE/LINKLENTH /LINK TWIST/ JOINT ANGLE or

JOINT DISTANCE

COG VECTOR

•

(for N links)

The LINKTYPE is given as 'R' for a revolute joint, and 'P' for a prismatic joint. The COG vector is given as ( $\rho_1$   $\rho_2$   $\rho_3$ ), and the inertia tensor J as ( $J_{11}$   $J_{12}$   $J_{13}$   $J_{22}$   $J_{23}$   $J_{31}$ ). In this way, the data for N links are given.

The data for the task of the manipulator, and the initial state are given in the file tASK.DAT in the following format:

INITIAL POSITION VECTOR

INITIAL VELOCITY VECTOR

A blank line or comment

FUNCTIONAL COORDINATES ACCELERATION

FUNCTIONAL COORDINATES ACCELERATION

•

•
FUNCTIONAL COORDINATES ACCELERATION

All these vectors are of the dimensions ( lXN ). The units for these values can be given in either MKS or CGS or FPS units. By default, MKS system of units is assumed.

## 4.4 RUNNING THE PROGRAM:

To start the program, the following monitor command is given

## .RUN ROBOTA

Now the control is passed to the top level CI and the flow of execution will now totally depend upon the commands given by the user.

#### CHAPTER V

#### RESULTS, CONCLUSIONS AND SUGGESTIONS

#### 5.1 RESULTS:

The program is run on a DEC - 10 computer using a sample input data. The input and output data, together with the transactions in a single session are listed in Appendix A 1. The listing illustrates some of the facilities provided by the Editor. Also appendend with it are the drive time, state time, and torque-rpm graphs. The output data obtained agrees with the actual values computed manually. The program listing is given in Appendix A 2.

#### 5.2 CONCLUSIONS:

As pointed out earlier, an algorithm which calculates the dynamic characteristics of a manipulator during the execution of a task clearly represents a useful means for the process of manipulater design. The algorithm chosen for this work forms and solves the dynamic model of the manipulator, and using incremental iterative displacement technique, computes the dynamic characteristics at a series of time instants. The dynamic model is based on the general theorem of Dynamics, and uses block matrix formalism to reduce the complex analytical equations into a compact form suitable for solving on a computer. The

inputs to the algorithm are the manipulator configuration (described by a set of parameters defining the geometric and inertial properties of the link), the initial state, and the manipulation task (in the form most suited for prescribing, according to the class of the functional movements). As the output, we obtain the time history of the generalised coordinates, velocities and drives, the P-N diagram, and the total energy consumed.

This algorithm is implemented in the increasingly popular Pascal language, to aid the designer in a fast analysis of the dynamic characteristics of a number of manipulator configuration. The P-N diagram can be used to give specifications for the actuator units. The total energy consumed for a particular task can be compared for different configurations and the configuration chosen accordingly. The various graphs can be viewed on a graphics terminal.

Any of the input parameters can be changed, by means of the Editor module and simulation done again. Further, the data may be given in any units. The program is command driven, i.e. the user types in commands for the action he desires. The command is decoded, and the action taken, by the Command Interpreter module. The Editor and

the Command Interpreter thus allow the user to analyse different configurations in an interactive manner. Further the structured development of the program allows new modules to be easily added.

### 5.3 SUGGESTIONS:

Improvements in the program can be made in two ways. One is an extension of the Dynamic module to include a manipulator with third and fourth class pairs, and a branched structure, and a manipulator with elastic segments. Secondly, other modules can be added to consider other problems in Robotics, like path synthesis, workspace synthesis actuator modelling, performance evaluation, etc. If this is done, coupled with the extensive use of computer graphics, then the gap between the designers and theoreticians can be bridged.

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## I - A XICHESSAN

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For complete help see the DUCUMENTATION.
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display joint accelerations
display Torque - Rom diagram
display energy consumption for
and output parameters display
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07

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   李巴门
  * HELP
                                                   The following is a summary of commands
                                                  THE STORY OF THE S
                                                                                                                                                                  type this text change the system of units, coordinates
                                                                                                                                                                  change the Manipulator change the Manipulator change the Manipulator
                                                                                                                                                                                                                                                                                                            configuration
state
task
                                                                                                                                                                  end édit
 *CONFIG
                                                   * Add, delete or change :
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                                                               Configuration parameter :
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\* Chapter :

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